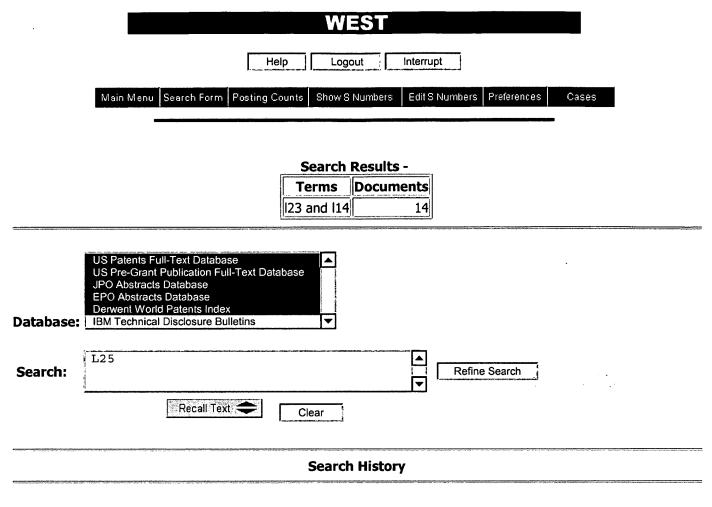
WEST
Freeform Search
US Patents Full-Text Database US Pre-Grant Publication Full-Text Database JPO Abstracts Database EPO Abstracts Database Derwent World Patents Index IBM Technical Disclosure Bulletins
126 not 129
10 Documents in Display Format: - Starting with Number 1
○ Hit List ● Hit Count ○ Side by Side ○ Image
Search Clear Help Logout Interrupt
Main Menu Show S Numbers Edit S Numbers Preferences Cases
Search History

DATE: Wednesday, March 06, 2002 Printable Copy Create Case

1 of 2

Set Name		<u>Hit Count</u>	Set Name result set		
DB=USPT,PGPB,JPAB,EPAB,DWPI; PLUR=YES; OP=ADJ					
<u>L30</u>	126 not 129	15	<u>L30</u>		
<u>L29</u>	L28 and (424/401.ccls. or cosmetic)	8	<u>L29</u>		
<u>L28</u>	118 and 16	26	<u>L28</u>		
<u>L27</u>	l23 and l18	1	<u>L27</u>		
<u>L26</u>	l6 and (424/401.ccls. or cosmetic)	23	<u>L26</u>		
<u>L25</u>	118 same 124	5	<u>L25</u>		
<u>L24</u>	ethylene glycol propylene glycol block copolymer or ethylene oxide propylene oxide block copolymer	693	<u>L24</u>		
<u>L23</u>	118 same 16	1	<u>L23</u>		
<u>L22</u>	l18 same (l3 and l14)	0	<u>L22</u>		
<u>L21</u>	l18 same (l3 and l6)	0	<u>L21</u>		
<u>L20</u>	I18 same I3	16	<u>L20</u>		
<u>L19</u>	L18 and I3	190	<u>L19</u>		
<u>L18</u>	oil-in-water or oil in water or o/w	22654	<u>L18</u>		
<u>L17</u>	l15 and (l14 or l6)	0	<u>L17</u>		
<u>L16</u>	L15 and I4	0	<u>L16</u>		
<u>L15</u>	l1 same l2	4	<u>L15</u>		
<u>L14</u>	113 same 112 same 19	5411	<u>L14</u>		
<u>L13</u>	L10 or I7	197142	<u>L13</u>		
<u>L12</u>	l11 or l8	125900	<u>L12</u>		
<u>L11</u>	propylene oxide	48307	<u>L11</u>		
<u>L10</u>	ethylene oxide	88299	<u>L10</u>		
<u>L9</u>	block copolymer	64428	<u>L9</u>		
<u>L8</u>	propylene glycol	94427	<u>L8</u>		
<u>L7</u>	ethylene glycol	135909	<u>L7</u>		
<u>L6</u>	poloxamer 124	64	<u>L6</u>		
<u>L5</u>	l1 same l2 same l3	0	<u>L5</u>		
<u>L4</u>	(ceteareth 20 or ceteareth-20)	407	<u>L4</u>		
<u>L3</u>	(ceteareth 20 or ceteareth-20)	407	<u>L3</u>		
<u>L2</u>	(arlacel 165 or simulsol 165)	333	<u>L2</u>		
<u>L1</u>	(pemulen tr1 or carbopol 1342)	244	<u>L1</u>		

END OF SEARCH HISTORY



DATE: Wednesday, March 06, 2002 Printable Copy Create Case

Set Name		Hit Count S	Set Name result set		
DB=USPT,PGPB,JPAB,EPAB,DWPI; PLUR=YES; OP=ADJ					
<u>L25</u>	I23 and I14	14	<u>L25</u>		
<u>L24</u>	L23 and (I12 or I11)	1	<u>L24</u>		
<u>L23</u>	I19 and (I9 or I10)	18	<u>L23</u>		
<u>L22</u>	l20 and (l11 or l12)	0	<u>L22</u>		
<u>L21</u>	I20 and (I9 or I10)	0	<u>L21</u>		
<u>L20</u>	L19 and I8	5.	<u>L20</u>		
<u>L19</u>	l17 and (l4 or l5)	75	<u>L19</u>		
<u>L18</u>	L17 and I3	0	<u>L18</u>		
<u>L17</u>	l1 and l2	1181	<u>L17</u>		
<u>L16</u>	I1 and I2 and (I4 or I5) and I6 and I7 and I8 and (I9 or I10) and (I11 or I12) and I13 and I14	0	<u>L16</u>		
<u>L15</u>	l1 and l2 and l3 and (l4 or l5) and l6 and l7 and l8 and (l9 or l10) and (l11 or l12) and l13 and l14	0,	<u>L15</u>		
<u>L14</u>	carboxyvinyl or carbomer	6106	<u>L14</u>		
<u>L13</u>	propylene glycol	94427	<u>L13</u>		
<u>L12</u>	poloxamer	2230	<u>L12</u>		
<u>L11</u>	poloxamer 124 or polyoxamer-124	64	<u>L11</u>		
<u>L10</u>	ceteareth	642	<u>L10</u>		
<u>L9</u>	ceteareth 20 or ceteareth-20	407	<u>L9</u>		
<u>L8</u>	arlacel 165 or simulsol 165	333	<u>L8</u>		
<u>L7</u>	peg 100 or peg-100 or polyethylene glycol 100	849	<u>L7</u>		
<u>L6</u>	glyceryl	16860	<u>L6</u>		
<u>L5</u>	pemulen tr1 or carbopol 1342	244	<u>L5</u>		
<u>L4</u>	acrylate/c10-c30 alkylacrylate or acrylate/c10-c30 alkyl acrylate	2	<u>L4</u>		
<u>L3</u>	nadifloxacin	9	<u>L3</u>		
<u>L2</u>	volatile silicone or cyclomethicone or cyclotetradimethylsiloxane or cyclopentadimethyl siloxane or cyclohexyldimethylsiloxane or dimethylsiloxane/methayalkyl siloxane or hexamethyl disiloxane or hexamethyldisiloxane or hexylheptamethyltrisiloxane or hexyl heptamethyltrisiloxane or octylheptamethyltrisiloxane or octyl heptamethyltrisiloxane or octyl heptamethyltrisiloxane or phenyl silicone	6931	<u>L2</u>		
<u>L1</u>	oil-in-water or o/w	22654	<u>L1</u>		

END OF SEARCH HISTORY

2 of 2

Trying 3106016892...Open

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PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

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=> fil reg

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

0.15

0.15

FULL ESTIMATED COST

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STRUCTURE FILE UPDATES: 4 MAR 2002 HIGHEST RN 398116-77-3 DICTIONARY FILE UPDATES: 4 MAR 2002 HIGHEST RN 398116-77-3

TSCA INFORMATION NOW CURRENT THROUGH July 7, 2001

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Calculated physical property data is now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details: http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

The P indicator for Preparations was not generated for all of the CAS Registry Numbers that were added to the H/Z/CA/CAplus files between 12/27/01 and 1/23/02. Use of the P indicator in online and SDI searches during this period, either directly appended to a CAS Registry Number or by qualifying an L-number with /P, may have yielded incomplete results. As of 1/23/02, the situation has been resolved. Also, note that searches conducted using the PREP role indicator were not affected.

Customers running searches and/or SDIs in the H/Z/CA/CAplus files incorporating CAS Registry Numbers with the P indicator between 12/27/01 and 1/23/02, are encouraged to re-run these strategies. Contact the CAS Help Desk at 1-800-848-6533 in North America or 1-614-447-3698, worldwide, or send an e-mail to help@cas.org for further assistance or to receive a credit for any duplicate searches.

```
=> s poloxamer 124/cn
L1 1 POLOXAMER 124/CN
```

=> d

L1 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS

RN 106392-12-5 REGISTRY

CN Oxirane, methyl-, polymer with oxirane, block (9CI) (CA INDEX NAME) OTHER NAMES:

CN Adeka 25R1

CN Adeka 25R2

CN Adeka L 61

CN Adeka Pluronic F 108

CN Antarox 17R4

CN Antarox 25R2

CN Antarox B 25

```
CN
           Antarox F 108
CN
           Antarox F 68
CN
           Antarox F 88
           Antarox F 88FL
CN
CN
           Antarox L 61
CN
           Antarox L 72
CN
           Antarox P 104
CN
           Antarox P 84
CN
           Antarox SC 138
CN
           Arco Polyol R 2633
                                                                                              to the stands of the stands of
CN
           Arcol E 351
CN
           B 053
CN
           BASF-L 101
CN
           Block polyethylene-polypropylene glycol
CN
           Block polyoxyethylene-polyoxypropylene
CN
           Breox BL 19-10
CN
           Cirrasol ALN-WS
           Crisvon Assistor SD 14
CN
           CRL 1005
CN
           CRL 1605
CN
CN
           CRL 8131
CN
           CRL 8142
CN
           D 500
CN
           D 500 (polyglycol)
           Daltocel F 460
CN
CN
           Detalan
CN
           DO 97
           Dowfax 30C05
CN
CN
           ED 56
CN
           Empilan P 7068
CN
           Emulgen PP 230
CN
           EP 3028
CN
           Epan 485
CN
           Epan 785
CN
           Epan U 108
CN
           Ethylene glycol-propylene glycol block copolymer
CN
           Ethylene oxide-propylene oxide block copolymer
CN
           Ethylene oxide-propylene oxide block copolymer dipropylene glycol ether
           Ethylene oxide-propylene oxide block polymer
CN
CN
           Ethylene oxide-propylene oxide copolymer, block
CN
           F 108
CN
           F 127
CN
           F 68
           Poloxamer 124
ADDITIONAL NAMES NOT AVAILABLE IN THIS FORMAT - Use FCN, FIDE, or ALL for
           11104-97-5, 163516-02-7, 124057-62-1, 121089-00-7, 96639-37-1,
96958-14-4,
           99040-06-9, 106138-19-6, 113441-83-1, 115742-90-0, 108688-61-5,
           108688-62-6, 37349-41-0, 70226-19-6, 72231-62-0, 77108-15-7, 80456-04-8,
           144638-32-4, 83589-65-5, 86904-45-2, 106899-85-8, 107498-07-7,
           108340-62-1, 188815-93-2, 211389-05-8, 355134-17-7
MF
            (C3 H6 O . C2 H4 O) \times
CI
           PMS, COM
PCT
           Polyether, Polyether formed
SR
LC
                                        ADISINSIGHT, ADISNEWS, AGRICOLA, ANABSTR, BIOBUSINESS,
           STN Files:
               BIOSIS, CA, CANCERLIT, CAPLUS, CASREACT, CEN, CHEMCATS, CHEMLIST, CIN,
                CSCHEM, DDFU, DIOGENES, DRUGNL, DRUGU, DRUGUPDATES, IPA, MEDLINE,
                PDLCOM*, PHAR, PIRA, PROMT, RTECS*, TOXCENTER, TOXLIT, USAN, USPAT2,
```

```
USPATFULL
         (*File contains numerically searchable property data)
     CM
     CRN 75-56-9
     CMF C3 H6 O
     CH3
     CM
          2
     CRN 75-21-8
     CMF C2 H4 O
            6144 REFERENCES IN FILE CA (1967 TO DATE)
             634 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA
            6174 REFERENCES IN FILE CAPLUS (1967 TO DATE)
=> s ceteareth 20/cn
             1 CETEARETH 20/CN
L2
=> d
    ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS
L2
   68439-49-6 REGISTRY *
* Use of this CAS Registry Number alone as a search term in other STN files
  result in incomplete search results. For additional information, enter HELP
  RN* at an online arrow prompt (=>).
    Alcohols, C16-18, ethoxylated (CA INDEX NAME)
OTHER NAMES:
CN
     Alfonic 1618-46
CN
     Alfonic 1618-60
     Alfonic 1618-62 ,
CN
CN
    Alfonic 1618-78
CN
    Alfonic 1618-80
CN
    B 2050-01A
CN
    Berol 07
CN
     Brij 68
CN
     C16-18 alcohols, ethoxylated
CN
     C16-18 ethoxylated alcs.
CN
     C16-18 fatty alcs., ethoxylated
CN
     Cemulsol DB
CN
     Ceteareth
CN
     Ceteareth 15
CN
    Ceteareth 20
CN
    Ceteareth 23
CN
    Ceteareth 25
```

```
CN
     Ceteareth 30
CN
     Cetomacrogol
CN
     Cetomacrogol 1000
CN
     Cetostearyl alcohol ethoxylated
CN
     Cremophor A 25
CN
     Empilan KM 11
CN
     Empilan KM 15
CN
     Empilan KM 20
CN
     Empilan KM 50
CN
     Emulgin ET 5L
CN
     Emulgin VP 3370
CN
     Ethal CSA 10
CN
     Ethal CSA 17
CN
     Ethoxylated alcohols, C16-18
CN
     Ethoxylated C16-18 alcs.
CN
     Ethoxylated cetostearyl alcohol
CN
     Ethoxylated cetostearyl alcohols
CN
     Eumulgin B 1
CN
     Eumulgin B 2
CN
     Eumulgin B 3
CN
     Eumulgin C 1000
CN
     Eumulgin C 1500
CN
     Eumulgin C 700
CN
     Fatty alcs., C16-18, ethoxylated
CN
     Hetoxol CS 25
CN
     Hetoxol CS 4
CN
     Hexotol CS 20
CN
     Imbentin AG 168/060
CN
     Imbentin AG 168/090
CN
     Imbentin AG 168/200
CN
     Imbentin AG 168S030
CN
     Imbentin AG 168S110
CN
     Imbentin AG 168S300
ADDITIONAL NAMES NOT AVAILABLE IN THIS FORMAT - Use FCN, FIDE, or ALL for
     DISPLAY
DEF
     This substance is identified by SDA Substance Name: C16-C18 alkyl
alcohol
     ethoxylate and SDA Reporting Number: 19-065-00. Consult SDA Substance
     Identification Procedure.
     12656-10-9, 12772-71-3, 12790-63-5, 8049-57-8, 54578-85-7, 54824-28-1,
DR
     55464-98-7, 59494-36-9, 60328-40-7, 64060-02-2, 64296-31-7, 64425-87-2,
     62887-14-3, 62887-15-4, 62887-16-5, 62887-17-6, 102381-15-7, 51877-40-8,
     37337-62-5, 68188-13-6, 85941-45-3, 39384-37-7, 39384-38-8, 39404-34-7,
     39404-35-8, 52732-71-5, 306966-46-1
MF
     Unspecified
CI
     MAN, CTS
LC
                  BIOSIS, CA, CAPLUS, CHEMCATS, CHEMLIST, CIN, CSCHEM, DDFU,
     STN Files:
       DRUGU, EMBASE, IFICDB, IFIPAT, IFIUDB, IPA, MSDS-OHS, PIRA, PROMT,
       RTECS*, TOXCENTER, TOXLIT, USPATFULL
         (*File contains numerically searchable property data)
     Other Sources:
                      DSL**, TSCA**
         (**Enter CHEMLIST File for up-to-date regulatory information)
*** STRUCTURE DIAGRAM IS NOT AVAILABLE ***
=> s nadifloxacin/c
NUMERIC VALUE NOT VALID 'NADIFLOXACIN'
L3
             0 NADIFLOXACIN/C
```

=> s nadifloxacin/cn

L4 1 NADIFLOXACIN/CN

=> d

L4 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2002 ACS

RN 124858-35-1 REGISTRY

CN 1H,5H-Benzo[ij]quinolizine-2-carboxylic acid, 9-fluoro-6,7-dihydro-8-(4-hydroxy-1-piperidinyl)-5-methyl-1-oxo- (9CI) (CA INDEX NAME)

OTHER CA INDEX NAMES:

CN 1H,5H-Benzo[ij]quinolizine-2-carboxylic acid, 9-fluoro-6,7-dihydro-8-(4-hydroxy-1-piperidinyl)-5-methyl-1-oxo-, (.+-.)-

OTHER NAMES:

CN Nadifloxacin

CN OPC 7251

DR 81962-84-7

MF C19 H21 F N2 O4

CI COM

SR CA

LC STN Files: ADISINSIGHT, BEILSTEIN*, BIOBUSINESS, BIOSIS, BIOTECHNO, CA, CAPLUS, CASREACT, CBNB, CIN, DDFU, DRUGNL, DRUGPAT, DRUGU, DRUGUPDATES, EMBASE, IPA, MEDLINE, MRCK*, PHAR, PROMT, SYNTHLINE, TOXCENTER, TOXLIT, USAN, USPATFULL

(*File contains numerically searchable property data)

$$Me$$
 HO_2C
 N
 N
 F

PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT

63 REFERENCES IN FILE CA (1967 TO DATE)

3 REFERENCES TO NON-SPECIFIC DERIVATIVES IN FILE CA

64 REFERENCES IN FILE CAPLUS (1967 TO DATE)

=> fil caplus uspatfull medline biosis embase kosmet COST IN U.S. DOLLARS SIN

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST 21.70 21.85

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```
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FILE 'BIOSIS' ENTERED AT 13:35:59 ON 06 MAR 2002
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COPYRIGHT (C) 2002 International Federation of the Societies of Cosmetics
Chemists
=> d his
     (FILE 'HOME' ENTERED AT 13:34:01 ON 06 MAR 2002)
     FILE 'REGISTRY' ENTERED AT 13:34:07 ON 06 MAR 2002
              1 S POLOXAMER 124/CN
L1
L2
              1 S CETEARETH 20/CN
L3
              0 S NADIFLOXACIN/C
L4
              1 S NADIFLOXACIN/CN
     FILE 'CAPLUS, USPATFULL, MEDLINE, BIOSIS, EMBASE, KOSMET' ENTERED AT
     13:35:59 ON 06 MAR 2002
=> s 11
'CN' IS NOT A VALID FIELD CODE
          8811 L1
=> s 12 or ceteareth 20
'CN' IS NOT A VALID FIELD CODE
           566 L2 OR CETEARETH 20
=> s 14 or nadifloxacin
'CN' IS NOT A VALID FIELD CODE
          224 L4 OR NADIFLOXACIN
=> s oil-in-water or o/w
'W' IS NOT A VALID FIELD CODE
        68752 OIL-IN-WATER OR O/W
1.8
=> s pemulen trl or pemulen tr 1 or carbopol 1342
           670 PEMULEN TR1 OR PEMULEN TR 1 OR CARBOPOL 1342
=> s arlacel 165 or simulsol 165
T-10
          447 ARLACEL 165 OR SIMULSOL 165
=> s propylene glycol
       119346 PROPYLENE GLYCOL
=> s carboxyvinyl or carbomer
L12
       7633 CARBOXYVINYL OR CARBOMER
```

```
(FILE 'HOME' ENTERED AT 13:34:01 ON 06 MAR 2002)
     FILE 'REGISTRY' ENTERED AT 13:34:07 ON 06 MAR 2002
              1 S POLOXAMER 124/CN
L1
              1 S CETEARETH 20/CN
L2
              0 S NADIFLOXACIN/C
L3
L4
              1 S NADIFLOXACIN/CN
     FILE 'CAPLUS, USPATFULL, MEDLINE, BIOSIS, EMBASE, KOSMET' ENTERED AT
     13:35:59 ON 06 MAR 2002
L5
           8811 S L1
L6
            566 S L2 OR CETEARETH 20
L7
            224 S L4 OR NADIFLOXACIN
L8
          68752 S OIL-IN-WATER OR O/W
L9
            670 S PEMULEN TR1 OR PEMULEN TR 1 OR CARBOPOL 1342
L10
            447 S ARLACEL 165 OR SIMULSOL 165
L11
         119346 S PROPYLENE GLYCOL
L12
           7633 S CARBOXYVINYL OR CARBOMER
=> s cyclomethicone or cyclotetradimethsiloxane or cyclopentadimethyl
siloxane or cyclohexyldimethylsiloxane or dimethylsiloxane/methyalkyl siloxne
or hexamethyl disiloxane or hexamethyldisiloxane or
hexylheptamethyltrisiloxane or hexyl heptamethyltrisiloxane or hexyl
heptamethyl trisiloxane
MISSING OPERATOR
MISSING OPERATOR
MISSING OPERATOR
MISSING OPERATOR
MISSING OPERATOR
MISSING OPERATOR
=> s volatile silicone
L13
         1645 VOLATILE SILICONE
=> d his
     (FILE 'HOME' ENTERED AT 13:34:01 ON 06 MAR 2002)
     FILE 'REGISTRY' ENTERED AT 13:34:07 ON 06 MAR 2002
L1
              1 S POLOXAMER 124/CN
L2
              1 S CETEARETH 20/CN
L3
              0 S NADIFLOXACIN/C
T.4
              1 S NADIFLOXACIN/CN
     FILE 'CAPLUS, USPATFULL, MEDLINE, BIOSIS, EMBASE, KOSMET' ENTERED AT
     13:35:59 ON 06 MAR 2002
L_5
           8811 S L1
L6
            566 S L2 OR CETEARETH 20
L7
            224 S L4 OR NADIFLOXACIN
L8
          68752 S OIL-IN-WATER OR O/W
            670 S PEMULEN TR1 OR PEMULEN TR 1 OR CARBOPOL 1342
L9
L10
            447 S ARLACEL 165 OR SIMULSOL 165
L11
         119346 S PROPYLENE GLYCOL
L12
           7633 S CARBOXYVINYL OR CARBOMER
L13
           1645 S VOLATILE SILICONE
=> s 18 and 113
          543 L8 AND L13
L14
```

=> s 114 and 112 and 111 and 110 and 19 and 16 L15 0 L14 AND L12 AND L11 AND L10 AND L9 AND L6

=> s 114 and 110 and 19 L16 5 L14 AND L10 AND L9

=> log y

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION FULL ESTIMATED COST 44.88 66.73

STN INTERNATIONAL LOGOFF AT 13:42:22 ON 06 MAR 2002